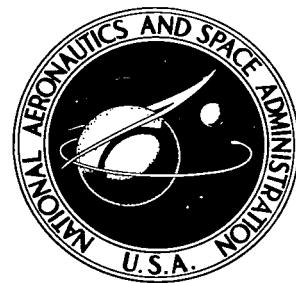


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SAVLOC, COMPUTER PROGRAM FOR
AUTOMATIC CONTROL AND ANALYSIS
OF X-RAY FLUORESCENCE EXPERIMENTS

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SAVLOC, COMPUTER PROGRAM FOR AUTOMATIC CONTROL AND ANALYSIS OF X-RAY FLUORESCENCE EXPERIMENTS

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SUMMARY

A program for a PDP-15 computer is presented which provides for control and analysis of trace element determinations by using X-ray fluorescence. The program simultaneously handles data accumulation for one sample and analysis of data from previous samples. Data accumulation consists of sample changing, timing, and data storage. Analysis requires the locating of peaks in X-ray spectra, determination of intensities of peaks, identification of origins of peaks, and determination of areal density of the element responsible for each peak. The program may be run in either a manual (supervised) mode or an automatic (unsupervised) mode.

INTRODUCTION

The recent development of high-resolution solid-state X-ray detectors has made possible the use of X-ray fluorescence as a tool for the detection of trace elements in a variety of applications. However, analysis of environmental samples by means of X-ray fluorescence generates energy spectra for which quantitative hand analysis is prohibitively tedious. In addition, the number of samples generated by any comprehensive monitoring program requires a speed of data acquisition and analysis which is possible only through the use of a computerized system. The present report describes the software designed for control and data analysis of such a system. This report should permit the reader to operate the system.

PRINCIPLES OF X-RAY FLUORESCENCE AND CHARACTER OF DATA

When properly stimulated, the atoms of any element emit X-rays whose energy (or wavelength) is strictly dependent upon the atomic number of the emitting atom. Since

these energies may be quickly and accurately measured, and since the X-ray energies of all elements are well known and tabulated, it is possible to identify, with little ambiguity, the element responsible for the emission. In addition, the number of X-rays emitted is proportional to the number of atoms stimulated, so that after a proper calibration, quantitative measurements of areal densities are possible.

The X-ray detectors commonly employed provide data in the form of analog electric pulses. Data acquisition consists of the digitizing and storing of these pulses, which arrive with random timing at the rate of several thousand per second. A schematic representation of the X-ray fluorescence facility is shown in figure 1. A typical X-ray spectrum is shown in figure 2. This particular spectrum resulted from excitation of a sample by using X-rays from a tungsten X-ray tube. It contains a peak (labelled Rh) which resulted from the presence in the primary X-ray beam of a small, known amount of rhodium, whose purpose was to provide a monitor of the number of photons incident on the sample.

Data analysis consists of determining from this spectrum the number of X-ray counts contained in each peak, identifying the origin of the peak, and converting the number to an areal density by using the expression

$$N(E) = N_i \left(\frac{d\sigma}{d\Omega} \right)_E \rho_x \Omega \epsilon(E) \quad (1)$$

where

$N(E)$	number of X-ray photons of energy E detected
N_i	number of X-ray photons incident on target
$(d\sigma/d\Omega)_E$	cross section for production of X-rays of energy E by exciting X-rays
ρ_x	areal density of atoms producing X-rays of energy E
Ω	solid angle subtended by X-ray detector
$\epsilon(E)$	efficiency of detector for detection of X-rays of energy E

In practice, only the quantity $N(E)$ is determined absolutely. The number N_i is assumed to have the form

$$N_i = K_{Rh} N_{Rh} \quad (2)$$

where N_{Rh} is the number of rhodium (monitor) X-rays detected, so that the areal density may be written

$$\rho_x = \frac{N(E)}{K_{Rh} N_{Rh} \left(\frac{d\sigma}{d\Omega} \right)_E \Omega \epsilon(E)} \quad (3)$$

The overall efficiency function

$$F(E) = \frac{1}{K_{Rh} \left(\frac{d\sigma}{d\Omega} \right)_E \Omega \epsilon(E)} \quad (4)$$

is then determined experimentally by using samples of known areal density of a number of elements. For elements other than those for which $F(E)$ is measured directly, an interpolation is carried out with the assumption that the function varies smoothly with energy. The areal density then, in terms of directly measurable quantities, becomes

$$\rho_x = N(E) \frac{F(E)}{N_{Rh}} \quad (5)$$

Analysis then requires only the determination of the intensity and energy of each line in the X-ray spectrum and the intensity of the line used for normalization N_{Rh} .

DATA ACQUISITION AND ANALYSIS

The entire X-ray fluorescence operation (sample changing, data acquisition, and data analysis) is carried out under the control of a PDP-15 computer. Communication between the computer and the X-ray fluorescence hardware is by means of a CAMAC system (ref. 1). The programmable interface between the computer and the CAMAC system has been described in detail elsewhere (ref. 2). The control panel of the PDP-15 computer, shown in figure 3, allows the input of a number of parameters as well as the selection of a number of different modes of operation, so that a sufficient degree of flexibility is provided to make the system useful for a variety of applications.

Data Acquisition

The two principal control tasks which the PDP-15 must perform are the changing of targets and the control (stop, start, and data transfer) of the 1024-channel analog-to-digital converter (ADC), through which all pulse-height data pass.

A typical data-taking cycle is initiated upon input of the last piece of sample identification data by means of the computer teletype unit. The ADC is turned on, and data transfers from the ADC to the computer memory take place autonomously by way of the CAMAC interface. During data acquisition the pulse-height spectrum is displayed on a video unit, which assures the user that all systems are functioning properly.

By means of push buttons on the computer control panel, data acquisition may be terminated at any time, either by pressing B4, which causes data to be recorded on magnetic tape, or by pushing B2, which ends the run without recording of the data. Data acquisition also terminates, with all accumulated data being recorded on tape, when the elapsed time reaches the preset value as entered (in tenths of seconds) on control panel thumb switch TH2.

Following termination and depending on the mode of operation selected, either another run is started automatically (T15 up) or a halt occurs (T15 down) until the operator chooses to continue by pressing B2. In either case sample changing may or may not be done, depending on the setting of rotary switch R1 on the control panel (sample is changed when R1=2).

The data recorded upon completion of a run include the run identification (a five-character file name and a three-digit file number), the time elapsed during data acquisition, and, of course, the pulse-height spectrum itself.

Because analog-to-digital conversion and data transfers take place at data-channel level, they proceed autonomously once initiated, so that the computer central processing unit is free at main-stream level to proceed with analysis of a previous data set if so instructed.

Data Analysis

Analysis of previous data sets may either be initiated by the operator (press B1) or take place automatically when the computer is operating in an automatic mode (T15 up). Analysis begins with the recalling of two previous runs from tape: one with the data of interest and one that is a background or blank run, which is subtracted to remove events resulting from scattering or impurities in the system or sample-supporting matrix.

When analysis is initiated by an operator, the program requests a file name and file number for each of these data sets. In the automatic mode the program searches for a data set with the same name as that previously analyzed, increases the file number to obtain the next data set, and uses the same background data set. Hence, it is always necessary that at least the first set be analyzed manually.

The first step in the analysis is the determination of the number of events N_{Rh} contained in the monitor (rhodium) peak for both the sample data and background data.

This is accomplished by locating the rhodium peak and then fitting that portion of the spectrum within 20 channels of it with a Gaussian peak plus linear background. The number of events within the Gaussian peak is then the N_{Rh} of equation (5). The amount of background in this region provides the normalization for subtraction of the blank spectrum.

It is assumed that since the blank spectrum results from an unloaded sample of the supporting matrix, it is identical in shape to that encountered in the loaded sample. Consequently, after normalization and smoothing over three channel segments, the two spectra are subtracted channel by channel to arrive at the spectrum to be analyzed.

The recalled spectrum (with background subtracted) then replaces the live display of data currently being accumulated on the video unit. Figure 4 shows the recalled spectrum derived from the data of figure 2. The number of events under each peak, that is, the number of detected X-ray photons of each energy, is determined by a least-squares fitting to each peak of a Gaussian distribution of adjustable height, width, and location. Initial estimates of peak locations may be supplied by the researcher by means of the light pen and video display (T11 down), or they may be located automatically by the program (T11 up). The automatic location procedure is described in appendix A.

Depending on the setting of toggle switches, an energy calibration may be supplied at this time (T16 down). The user may supply through the teletype console the energies of the first two peaks selected with the light pen. For operation in the automatic mode (T16 and T15 up) the program retains the energy calibration from the previous run. Hence, again at least one run must be done manually. The researcher is also asked to supply through the teletype console (unless the computer is operating in the automatic mode) an estimate of the full width at one-half of the maximum (FWHM) of a typical peak.

The fitting routine does each peak in turn from the lowest to the highest energy. When two peaks are found to lie within four times the estimated FWHM, they are fitted simultaneously. Up to five peaks may be treated as a multiplet and done simultaneously. A more detailed description of the fitting procedure is given in appendix B.

Following analysis of each line for energy and intensity, an identification is made by comparing the measured energy with a table of energies and atomic symbols. The elements which are included in this table are listed in table I. If a peak is identified as a K_α line, a check is made to determine whether the K_β line can also be observed. If a line of the proper energy is located, the expected intensity of the K_β line (determined from the intensity of the K_α line) is subtracted from the K_β candidate. The remainder of the peak is tested for other identification. No peak may be identified as an L series X-ray unless both the L_α and L_β lines are observed.

The experimental intensity is converted to an areal density by using the stored efficiency function $F(E)$. This function is assumed (for K series X-rays) to have the form

$$\log \left[\frac{1}{F(E)} \right] = \sum_{i=1}^7 \frac{P(i)}{E^{(i-1)}} \quad (6)$$

The parameters $P(i)$ are determined by measuring the X-ray yield for a number of samples (approximately 10 to 15) of known areal density and performing a least-squares fit of the function given in equation (6) to that data. The parameters $P(i)$ vary with any change in detector or source geometry and must be redetermined each time such a change occurs. A typical set of parameters $P(i)$ is listed in table II. For heavier elements, which are measured by means of D series X-rays, tabulated efficiencies are used for those elements which the program can identify.

The estimates of errors in the computed areal density are based on the number of counts in each peak, given by

$$N(E) = I(E)\sigma(E)\sqrt{\pi} \quad (7)$$

where I and σ (as defined in appendix B) are the height and width, respectively, of the Gaussian distribution derived from the data. The areal density, given by equation (5), then becomes

$$\rho_X = \frac{I(E)\sigma(E)\sqrt{\pi}F(E)}{N_{Rh}} \quad (8)$$

Therefore, assuming the error in N_{Rh} to be small gives

$$\frac{\delta\rho_X}{\rho_X} = \left| \frac{\delta I(E)}{I(E)} \right| + \left| \frac{\delta\sigma(E)}{\sigma(E)} \right| + \left| \frac{\delta F(E)}{F(E)} \right| \quad (9)$$

or

$$\frac{\delta\rho_X}{\rho_X} = \left| \frac{\delta I(E)}{I(E)} \right| + \left| \frac{\delta\sigma(E)}{\sigma(E)} \right| + \left| \frac{1}{F(E)} \frac{dF(E)}{dE} \delta E \right| \quad (10)$$

The errors in each of the quantities $I(E)$, $\sigma(E)$, and E are derived in the usual way from the residuals of the least-squares fitting process.

Computed areal densities may be modified by an arbitrary multiplicative factor G , which may be entered by means of thumb wheel TH1. The value of G may be between

0 and 9.9999, and its effect is included in all areal densities and errors listed as output. The principal purpose of this factor is to permit the user to convert to a convenient system of units if so desired. If G is set equal to 1, the results are in units of micrograms per square inch.

Output may consist of either one or two pages of teletype output as well as a record of the results on magnetic tape. These options are selected through the control panel toggle switches (p. 1, T14 up; p. 2, T12 up; and tape, T13 up). Samples of the two pages of teletype output are shown in figures 5 and 6. More detailed information on the results of the least-squares fitting process may be obtained by replying "T" (true) when the program inquires, through the teletype console, "Do you wish extra output? Output=." A sample of this augmented output is shown in figure 7. It includes the height, width, and location, as well as the standard deviation of each, for each line fitted by the program.

A flow chart for the entire program is shown in figure 8. A guide to the loading of the program into the PDP-15 is given in appendix C. Appendix D contains the details of the overlay structure necessary in order to fit the program into the available core. Appendix E contains a listing of the FORTRAN sections of the program.

CONCLUDING REMARKS

The program described in this report permits trace element analysis of environmental samples at a rate which would make meaningful monitoring surveys feasible. Specifically, data analysis requires between 3 and 12 minutes, depending on the complexity of the spectrum, while data acquisition requires 5 to 30 minutes, depending on the primary radiation source and associated hardware employed. The program works reliably in an automatic mode; it changes samples, takes data, locates peaks, measures intensities, and identifies elements. As a result samples may be processed without an operator in attendance, which makes around-the-clock analysis possible. In addition, when operated under the supervision of the user, the program is capable of treating a wide variety of nonroutine samples, such as those containing elements not automatically recognized and those containing a number of heavy elements and thus having extremely complex spectra. For these, the program permits the extraction of meaningful data with a minimum amount of effort on the part of the user.

Lewis Research Center,
National Aeronautics and Space Administration,
Cleveland, Ohio, December 2, 1976,
506-25.

APPENDIX A

AUTOMATIC PEAK LOCATING PROCEDURE

If the operator so desires, initial estimates of peak locations are determined by the program, using a method described by Mariscotti (ref. 3). The method is based on the construction of smoothed second differences of the experimentally measured energy spectrum. The smoothed second difference is defined by

$$S_i(z, m) = \underbrace{\sum_{j=i-m}^{i+m} \dots \sum_{h=l-m}^{l+m}}_z (N_{l+1} - 2N_l + N_{l-1}) \quad (A1)$$

or equivalently

$$S_i(z, m) = \sum_{\text{all } j} C_{ij}(z, m) N_j \quad (A2)$$

where the C_{ij} are weighting factors derived by Mariscotti, given by

$$C_{ij}(z, m) = \sum_{l=i-m}^{i+m} C_{lj}(z-1, m) \quad (A3)$$

and

$$C_{ij}(0, m) = \begin{cases} 0 & \text{if } |j - i| \geq 2 \\ 1 & \text{if } |j - i| = 1 \\ -2 & \text{if } j = i \end{cases} \quad (A4)$$

The values of z and m used here are those found by Mariscotti to be most suitable, namely,

$$z = 5$$

$$m = \frac{(0.6\Gamma - 1)}{2}$$

The resulting function S then vanishes for a linear spectrum and is similar to the second derivative of a Gaussian function if N_j is a Gaussian function. The program assumes that a peak exists in channel i if three conditions are present: First,

$$|S_i| > 1.6 |\delta S_i| \quad (A5)$$

where δS_i is the standard deviation in the function S_i , approximated by Mariscotti as

$$\delta S_i = \left(\sum C_{ij}^2 \right)^{1/2} \delta N_i$$

where δN_i is the standard deviation in the data point N_i . The program assumes

$$\delta N_i = (N_i + B_i)^{1/2}$$

where B_i is the amount of background previously subtracted from channel i of the spectrum. Second,

$$S_i < 0 \quad (A6)$$

Third,

$$S_{i-1} < |S_i| > S_{i+1} \quad (A7)$$

The code does not reject any peak on the basis of its width as suggested by Mariscotti.

APPENDIX B

LEAST-SQUARES FITTING

Once initial estimates of peak locations and widths have been obtained, a least-squares fitting is carried out by using the Gaussian function

$$f(x) = I \exp \left[\frac{(x - x_0)^2}{\sigma^2} \right] \quad (B1)$$

to represent that portion of the spectrum being treated. In the event that several peaks fall within 4FWHM of each other, they are fitted simultaneously as the sum of as many as five Gaussian distributions, each of which has independently variable width.

For initial estimates, x_0 is taken to be the channel i either found by the automatic search routine described in appendix B or supplied through the light pen. The initial estimate of the intensity I is taken to be the number of counts in that i^{th} channel, and σ is calculated from the FWHM input (for a Gaussian distribution, $\sigma = \text{FWHM}/1.665$).

Fitting is carried out by standard methods. In the event a satisfactory fit cannot be obtained within a specified number of iterations, the peak of highest energy is dropped from the multiplet, and the fitting process is restarted for the remainder of the peaks. In the event that a single peak cannot be satisfactorily fitted, a message to that effect is included on output page 1, and the peak is assigned intensity zero and a location equal to the starting channel number.

APPENDIX C

LOADING THE PROGRAM

As presented in this report the program (entitled SAVLOC) is too large to be loaded directly into a 16K PDP-15 computer. As a result it can only be used as a series of overlays, constructed by using the routine CHAIN (ref. 4). A listing of the execute file, titled SAVLOC, is given in appendix D; the listing includes the resident code as well as the structure of each of the links and a machine constructed map of the file.

At execution time the user must assign handlers to three slots in the device assignment table (dat): dat slot 1, for writing of raw data; dat slot 2, for reading of raw data; and dat slot 3, for writing of results of analysis (fig. 6). Dat slots 1 and 2 are normally assigned the same magnetic tape (dectape) unit, and dat slot 3 is assigned a second dectape unit.

An example of the entire loading sequence is as follows:

KMS15 V5C
\$A DTE6 1,2/DTE7 3
\$E SAVLOC

EXECUTE V4A
FNAME(A) =

where the underlined expressions are output from the computer.

APPENDIX D

OVERLAY STRUCTURE AND MEMORY MAP FOR EXECUTE FILE SAVLOC

\CHAIN

CHAIN V5B

```
NAME XCT FILE
>SAVLOC
LIST OPTIONS & PARAMETERS
>PGR,16K,SZ
DEFINE RESIDENT CODE
>SIMUL2,CAMADC,IPK3,PICKER
DESCRIBE LINKS & STRUCTURE
>LK1=PEDE,RITE
>LK2=FITALL,FITALI,GAUSS
>LK3=ECAL1,ECAL2
>LK4=REGOUT,XKYLD
>LK5=LAREL,SOFT
>LK6=S,SCAL1,LOCUM
>LK1:LK2:LK3:LK4:LK5:LK6:PEED:MAXER:SET1:SUBTR:OUTER:
-MULT:SIT2:EPPO:ENERG:IDENT
>
```

LINK TABLE
37225-37636 00412

RESIDENT CODE

SIMUL2	36527-37224	00476
CAMADC	35731-36526	00576
IPK3	34747-35730	00771
PICKER	34305-34737	00433
RELAON	34226-34304	00057
RTIMER	34146-34225	00260
RCV16	34126-34145	00000
SEND16	34113-34125	00013
CAMAC.	33631-34112	00062
PANEL	33477-33630	00130
CROSS	33347-33476	00130
HISTO1	33226-33346	00121
PENSW	33200-33225	00026
VPVECT	33014-33177	00164
VIDEO.	32515-33013	00277
RTIME	32470-32514	00025
CLOCK.	32372-32467	00076
RTEST	32350-32371	00022
GETBLK	32267-32347	00061
IPOOL	32224-32266	00043
ZERO	32200-32223	00024
.SS	32130-32177	00050
FLOAT	32117-32127	00011
FLOAT2	32271-32116	00226
.BR	32224-32070	00045
FIOPS	31302-32023	00522
INTEAE	31171-31301	00111
PELEAE	30257-31170	01112

OTSEP 27712-27777 00066
.DA 30021-30056 00036
DISPL 30016-30020 00003
ERP 30012-30015 00004
STUFF 27464-27711 00226
MUP 27445-27463 00017
FINOUT 27345-27444 00100

LINK -- LK1
REDE 27277-27344 00046
PITE 27135-27276 00142
RELAOF 27061-27134 00054
BCDEXT 27012-27060 00047
INPUT1 26723-27011 00067
XREAD 26497-26702 00314
PACKER 24306-26406 00101
MOD 26265-26305 00021
BCDIO 23630-26264 02435
BINIO 23403-23627 00225
FILE 23124-23402 00257
.CR 23105-23123 00017

LINK -- LK2
FITALL 26136-27344 01207
FITALI 26005-26135 00131
GAUSS 25277-26004 00506
MATINV 24736-25276 00341
GENBLK 24663-24735 00053
LOC 24654-24662 00007
ABS 24635-24653 00017
RMNMX 24511-24634 00124
.RC 24415-24510 00074
EXP 24377-24414 00016
.EF 24272-24376 00105
.EC 24220-24271 00052
.RDA 24130-24217 00070
FITALC 24127-24127 00001

LINK -- LK3
ECAL1 27001-27344 00344
ECAL2 26446-27000 00333
INPUT1 26357-26445 00067
XREAD 26043-26356 00314
PACKER 25742-26042 00101
BCDIO 23305-25741 02435
.CR 23266-23304 00017

LINK -- LK4
REGOUT 26470-27344 00655
XKYLD 26343-26467 00125
.BC 26247-26342 00074
.BE 26237-26246 00010
.PR 26201-26236 00036
SQRT 26066-26200 00113
.EE 25703-26065 00163
.EF 25576-25702 00105
.EC 25524-25575 00052
.RDA 25434-25523 00070

BCDIO 22777-25433 02435
.CB 22760-22776 00017

LINK -- LK5
LABEL 24757-27344 02366
SOPT 24621-24756 00136
ABS 24602-24620 00017

LINK -- LK6
S 27117-27344 00226
SCALI 24651-27116 00246
LOCUM 24730-26650 01721
IFIX 24712-24727 00016
ARS 24673-24711 00017
LIMIT 24650-24672 00023
.BC 24554-24647 00074
SQRT 24441-24553 00113
DIFF 24220-24440 00221

LINK -- REED
PEED 26775-27344 00350
BCDEXT 26726-26774 00047
INPUT1 26637-26725 00067
XPEAD 26323-26636 00314
PACKER 24222-26322 00101
MOD 26201-26221 00021
BCDIO 23544-26200 02435
BINIO 23317-23543 00225
FILE 23040-23316 00257
.CR 23021-23037 00017

LINK -- MAXER
MAXER 27275-27344 00050

LINK -- SET1
SET1 27003-27344 00342

LINK -- SUBTP
SUBTP 26527-27344 00616
INPUT1 26440-26526 00267
XPEAD 26124-26437 00314
PACKER 26023-26123 00101
BCDIO 23366-26022 02435
.CR 23347-23365 00017

LINK -- OUTER
OUTER 27041-27344 00304
BCDIO 24404-27040 02435
.CB 24365-24403 00017

LINK -- MULT
MULT 26665-27344 00460
IFIX 26647-26664 00016

LINK -- SIT2
SIT2 26032-27344 01313
BCDIO 23375-26031 02435
.CR 23356-23374 00017

LINK -- ERRO
ERRO 27033-27344 00312
BCDIO 24376-27032 02435
.CB 24357-24375 00017

LINK -- ENRG
ENERG 26500-27344 00645

LINK -- IDENT
IDENT 26066-27344 01257
BCDEXT 26017-26065 00047
PACKER 25716-26014 00101
ARS 25677-25715 00017
MOD 25656-25676 00021
BCDIO 23221-25655 02435
FILE 22742-23220 00257
.CR 22723-22741 00017

BLANK COMMON
.XX 12723-22722 10000

CORE REQ'D
12723-37636 24714

KMS15 V5C
\$

APPENDIX E

SOURCE LISTING

```
C.....PROGRAM TO SIMULTANEOUSLY TAKE AND ANALYZE DATA
      INTEGER DATAD(4096), GATEX(64,7), ORGN(8), DATUM(1024)
      REAL FNAME(2), LT, LTI
      LOGICAL ON, EXIST, BUTTNS, RITEST
      EXTERNAL ADCON
      COMMON DATAD
      EQUIVALENCE (GATEX(1,1),DATAD(3573)),(DATAD(1),DATUM(1))
      DATA ND, NLAST/1024,1/,
     1   FNAME(1), FNAME(2)/5HSPECT,1H /
      CALL CLKON
      CALL PSTART
2     CALL REDE(FNAME, NEXT, NLAST)
3     I4=IROTOR(4)
     I3=IROTOR(3)
     I2=IROTOR(2)
     EXIST=.TRUE.
     NY=IROTOR(4)+6
     NF=256*I2-255
     NX=I3+7
     CALL ZERO(DATUM)
     CALL CLKON
     CALL VPSTOP
     CALL HISTO1(DATUM, NF, NX, NY, NSAW, 1)
     CALL VPSTRT
     CALL ECAMCK
     CALL RTCOFF
104    IF(ON) CALL ADCOFF
     MODE=0
     IRT=ITHUMB(2)
     ILT=0
     IF(NEXT.GT.NLAST) GO TO 102
     CALL ADCSTR(DATAD, GATEX, ORGN, ILT, ON, MODE)
     CALL RTCSTR(IRT, ADCOFF)
     CALL RTIME(TS)
     CALL LAMPON(1)
102    CALL BUTTNS(IR)
     IF(IR.EQ.1) GO TO 101
     IF(IR.EQ.2) GO TO 501
     IF(.NOT.ON) GO TO 96
102    IF(IR.EQ.4) GO TO 98
     CALL TOGGLS(IT)
     IF(BITEST(IT,15).AND.EXIST) GO TO 101
     GO TO 100
101    CALL VPSTOP
     CALL IPK2(EXIST)
     CALL VPSTOP
     CALL HISTO1(DATUM, NF, NX, NY, NSAW, 1)
     CALL VPSTRT
99     IF(ON) GO TO 102
96     IF(NEXT.GT.NLAST) GO TO 102
98     CALL RTCOFF
     CALL ADCOFF
     CALL LAMPOF(1)
```

```

CALL TOGGLS(IT)
IF(IROTOR(1).EQ.2) CALL RELAON(3)
CALL RITE(FNAME,NEXT,TS,IRT,ILT,ND)
NEXT=NEXT+1
CALL TOGGLS(IT)
IF(BITEST(IT,15)) GO TO 3
CALL LAMPON(2)
500 CALL BUTTNS(IB)
IF(IB.NE.2) GO TO 500
502 CALL LAMPOF(2)
501 CALL TOGGLS(IT)
IF(BITEST(IT,17)) GO TO 2
GO TO 3
END

SUBROUTINE IPK2(EXIST)
C.....SUBROUTINE TO DO PEAK ANALYSIS AFTER LIGHT PEN SELECTION,
C.....FOR USE WITH SIMUL2, TO TAKE DATA SIMULTANEOUSLY.
C.....USED TO BUILD XCT FILE
INTEGER LIST(520),DATA(1024),BKGND(1024),GATE(5,30),CH0(30),
1 CHMAX,DATAD(4096),BEXT,IR(5)
REAL FNAME(2),PARAM(15),DP(15),POINT(100),BNAME(2),XCH(2),
1 ELE(30),ENORM(30),BNORMX(30),DR(30)
LOGICAL EXIST,DOUBLE,UNDOUR,TRIPLE,UNTRIP,QUAD,UNQUAD,QUIN,
1 UNQUIN,BUTTNS,PENSW,OUTPT,BITEST
EXTERNAL GAUSS
COMMON/DISPL/NF,NX,NY
COMMON/ERR/PHI2,IMIN,ITELL
COMMON/STUFF/GATE
COMMON/MUP/DOUBLE,UNDOUR,TRIPLE,UNTRIP,QUAD,UNQUAD,QUIN,UNQUIN,
1 JPEAK,GAM,IDUB,ITRIP,IQUAD,IQUIN
COMMON/FINOUT/ELE,A1,A2
COMMON DATAD
EQUIVALENCE (DATA(1),DATAD(1025)),(BKGND(1),DATAD(2049)),
1 (LIST(1),DATAD(3073)),(CH0(1),DATAD(4021)),
2 (POINT(1),DATAD(3593)),(ENORM(1),DATAD(3800)),
3 (BNORMX(1),DATAD(3861)),(DR(1),DATAD(3921)),
4 (IR(1),DATAD(4051)),(PARAM(1),DATAD(4056))
CALL LAMPON(3)
CALL REED(1,FNAME,NEXT,EXIST)
IF(EXIST) GO TO 3
CALL LAMPOF(3)
RETURN
3 NEXT=NEXT+1
2 CALL MAXER(DATA,CHMAX)
CALL SETI(CHMAX,DATA,POINT)
ITELL=1
CALL FITALL(GAUSS,POINT,PARAM,DP,Q,LIST)
CALL OUTER(PARAM,LIST,ANORM,ARNORM,CHMAX)
CALL REED(2,RNAME,BEXT,EXIST)
CALL MAXER(BKGND,CHMAX)
CALL SETI(CHMAX,BKGND,POINT)
CALL FITALL(GAUSS,POINT,PARAM,DP,Q,LIST)
ITELL=0
CALL SUBTR(ANORM,ARNORM,CHMAX,OUTPT,GAM)
CALL PICKER
CALL TOGGLS(IT)
IF(BITEST(IT,10)) CALL UPSTOP

```

```

        IF(BITEST(IT,11)) CALL LOCUM
        CALL LAMPOF(7)
        IF(BITEST(IT,16)) GO TO 87
        DO 88 J=1,2
        CALL ECALI(J)
        CALL FITALL(GAUSS,POINT,PARAM,DP,Q,LIST)
        XCH(J)=FLOAT(CH0(J))+PARAM(2)-4.
88      CONTINUE
87      CALL ECAL2(XCH,A1,A2)
C.....BEGIN FITTING PEAKS
80      CALL LAMPOF(7)
        DO 89 J=3,JPEAK
        CALL MULT(IMAX,IMIN,J)
        CALL SIT2(IMAX,IMIN,J,ANORM)
        CALL FITALL(GAUSS,POINT,PARAM,DP,Q,LIST)
        NPARAM=LIST(6)
        DO 85 I=1,NPARAM,3
        IF(PARAM(I+1).GT.FLOAT(IMAX-IMIN+5)) LIST(4)=7
        IF(PARAM(I+1).LT.-5.) LIST(4)=8
85      CONTINUE
        IF(LIST(4).EQ.0) GO TO 86
        CALL ERRO(J,IMIN,DP)
        IF(.NOT.DOUBLE) GO TO 86
        GO TO 81
86      CALL ENERG(IMIN,A1,A2,J,DP)
        CALL LABEL(J,JPEAK)
        CALL REGOUT(OUTPT,ANORM,IMIN,DP,Q,J)
89      CONTINUE
        CALL IDENT(ELE,ENORM,BNORMX,DB,JPEAK,FNAME,NEXT)
        CALL LAMPOF(3)
        CALL VPSTRT
        RETURN
        END

```

```

SUBROUTINE PICKER
INTEGER GATE(5,30),CH0(30),DATA(1024),DATAD(4096),OFF,X,Y,
I      CHMAX
LOGICAL DOUBLE,UNDOUB,TRIPLE,UNTRIP,QUAD,UNQUAD,QUIN,UNQUIN,
I      BUTINS,PENSW,BITEST
COMMON/STUFF/GATE
COMMON/MUP/DOUBLE,UNDOUB,TRIPLE,UNTRIP,QUAD,UNQUAD,QUIN,
I      UNQUIN,JPEAK,GAM,IDUB,ITRIP,IQUAD,IQUIN
COMMON/DISPL/NF,NX,NY
COMMON DATAD
EQUIVALENCE (DATA(1),DATAD(1025)),(CH0(1),DATAD(4021))
DATA OFF/-1/
100    FORMAT(1X,2HOK)
        ICHK=0
6       ISI=NX
        IS2=NY
        IS3=NF
        CALL SCALI(NX,NY,NF)
        CALL TOGLLS(IT)
        IF(ICHK.NE.1) GO TO 66
        IF((ISI.NE.NX).OR.(IS2.NE.NY).OR.(IS3.NE.NF)) GO TO 66
        GO TO 65
66      CALL VPSTOP
        CALL HISTO1(DATA,NF,NX,NY,NHIT,1)

```

```

      CALL VPVECT(GATE,30,IHIT,2)
      CALL VPSTRT
65     IF(BITESIT(IT,16)) RETURN
      IF(ICCHK.EQ.1) GO TO 12
      IF(ICCHK.EQ.-1) GO TO 13
7       CALL ZERO(GATE)
      CALL ZERO(CH0)
      JPEAK=0
      CALL CROSS(X,Y,OFF,3)
8       JPEAK=JPEAK+1
12     IF(PENSW(.TRUE.)) GO TO 9
      ICHK=1
      GO TO 6
9       GATE(1,JPEAK)=NHIT
      CH0(JPEAK)=NHIT
      GATE(2,JPEAK)=DATA(NHIT)
      GATE(3,JPEAK)=0
      GATE(4,JPEAK)=128
      GATE(5,JPEAK)=-10
      ICHK=-1
      GO TO 6
13    CALL LAMPON(7)
10     CALL BUTINS(IB)
      IF(IR.EQ.128) GO TO 11
      IF(IB.EQ.32) GO TO 14
      IF(IB.EQ.64) RETURN
      GO TO 10
14    JPEAK=JPEAK-1
11     CALL LAMPOF(7)
      GO TO 8
      END

      SUBROUTINE REDE(FNAME,NEXT,NLAST)
      REAL FNAME(2)
      CALL INPUTI(FNAME(1),5HFNAME,6)
      CALL INPUTI(NEXT,4HNEXT,1)
      CALL INPUTI(NLAST,5HNLAST,1)
      RETURN
      END

      FUNCTION S(ICHLN)
      INTEGER DATA(1024),DATAD(4096)
      REAL C(36),B(36)
      COMMON/DIFF/C,B,M
      COMMON DATAD
      EQUIVALENCE (DATA(1),DATAD(1025))
      MI=5*M*2
      X=C(1)*FLOAT(DATA(ICHLN))
      DO 41 J=1,MI
      I1=ICHLN+J
      I2=ICHLN-J
      J1=J+1
      IF(I1.GT.1024) GO TO 42
      IF(I2.LT.0) GO TO 42
      X=X+C(J1)*FLOAT(DATA(I1))+C(J1)*FLOAT(DATA(I2))
      GO TO 41

```

```

42      X=0
41      CONTINUE
S=X
RETURN

SUBROUTINE RITE(FNAME,NEXT,TS,IRT,ILT,ND)
REAL FNAME(2)
INTEGER DATAD(4096)
COMMON DATAD
CALL BCDEXT(FNAME,NEXT)
CALL ENTER(2,FNAME)
WRITE(2) TS,IRT,ILT,ND,(DATAD(I),I=1,ND)
CALL RELAOF(3)
WRITE(6,106) FNAME,TS,IRT,ILT
106   FORMAT(IX,2A5,5X,F10.5,I6,3X,I6)
CALL CLOSE(2)
RETURN
END

```

```

SUBROUTINE GAUSS(R,X,P,DP,V,KK,J)
C.....SUBROUTINE FOR US EWITH FITALL TO FIT UP TO 4
C.....GAUSSIAN PAKS WITH OR WITHOUT BACKGROUND
REAL X(100),P(15),DP(15)
INTEGER BKGND(1024),DATA(1024),DATAD(4096)
COMMON/ERR/PHI2,ILOWER,ITELL
COMMON DATAD
EQUIVALENCE (DATA(1),DATAD(1025)),(BKGND(1),DATAD(2049))
NP=KK/3
IF(KK-3*NP.EQ.0) GO TO 1
R=X(J)-P(1)*FLOAT(J)-P(2)
DP(1)=-FLOAT(J)
DP(2)=-1.
N=3
GO TO 3
1  N=1
R=X(J)
3  DO 2 I=N,KK,3
TP=P(I)*EXP(-(FLOAT(J)-P(I+1))**2/P(I+2)**2)
R=R-TP
DP(I)=-TP/P(I)
DP(I+1)=-TP*2.*((FLOAT(J)-P(I+1))/P(I+2)**2
DP(I+2)=DP(I+1)*(FLOAT(J)-P(I+1))/P(I+2)
2  CONTINUE
IERR=ILOWER+J-3
IF(ITELL.EQ.0) V=(FLOAT(BKGND(IERR)+1))
IF(ITELL.EQ.1) V=1.
RETURN
END

```

```

SUBROUTINE ECALI(J)
C.....SUBROUTINE CALL ED BY IPK2 TO DETERMINE ENERGY CALIBRATION.
INTEGER CH0(30),LIST(520),DATAD(4096),DATA(1024)
LOGICAL BITEST
REAL PARAM(15),DP(15),POINT(100),X(2)
COMMON DATAD
COMMON/ERR/PHI2,IMIN,ITELL

```

```

EQUIVALENCE (DATA(1),DATAD(1025)),(LIST(1),DATAD(3073)),
1      (CH0(1),DATAD(4021)),(POINT(1),DATAD(3593)),(PARAM(1)
1,DATAD(4056))
CALL TOGLLS(IT)
IF(BITEST(IT,16)) RETURN
LIST(1)=7
LIST(2)=3
LIST(3)=15
LIST(4)=2
LIST(5)=3
LIST(6)=3
IK=CH0(J)
PARAM(1)=FLOAT(DATA(IK))
PARAM(2)=4.
PARAM(3)=4.5/1.665
IMIN=CH0(J)-3
DO 3 I=1,7
IC=CH0(J)-4+I
POINT(I)=FLOAT(DATA(IC))
CONTINUE
RETURN
END
3

```

```

C.....SUBROUTINE ECAL2(X,A1,A2)
SUBROUTINE TO DO ENERGY CALIBRAION WITH IPK2
REAL X(2)
LOGICAL BITEST
CALL TOGLLS(IT)
IF(BITEST(IT,16)) GO TO 99
WRITE(6,100)
CALL INPUT1(E1,2HE1,2)
CALL INPUT1(E2,2HE2,2)
A1=(E1-E2)/(X(1)-X(2))
A2=E1-A1*X(1)
99   G=FLOAT(ITHUMB(1))/10000.
WRITE(6,103) G
103  FORMAT(1X,2HG=,F6.4)
IF(BITEST(IT,14)) WRITE(6,101)
IF(BITEST(IT,14)) WRITE(6,102)
RETURN
101  FORMAT(///,6H PEAK,9H CHANNEL,12H STRENGTH,
1      10H ENERGY,16H CONCENTRATION,10H ERROR,
2      8H ELEM )
102  FORMAT(39X,14H(UGRAMS/IN**2))
100  FORMAT(1X,25HENERGIES FOR CALIBRATION:)
END

```

```

SUBROUTINE REGOUT(OUTPT,ANORM,IMIN,DP,Q,J)
INTEGER LIST(520),DATAD(4096)
REAL PARAM(15),ENORM(30),ELE(30),DP(15),BNORMX(30),DB(30)
LOGICAL OUTPT,BITEST
COMMON DATAD
COMMON/FINOUT/ELE,A1,A2
EQUIVALENCE (LIST(1),DATAD(3073)),(ENORM(1),DATAD(3800)),
1      (BNORMX(1),DATAD(3861)),(DB(1),DATAD(3921)),
2      (PARAM(1),DATAD(4056))
NPARAM=LIST(6)

```

```

DO 92 I=1,NPARAM
DP(I)=SQRT((Q)*DP(I))
92 CONTINUE
DO 91 I=1,NPARAM,3
IP2=I+1
ANORMX=PARAM(I)*PARAM(I+2)*1.7725/ANORM
BNORMX(J)=ANORMX*XYLD(ENORM(J))*FLOAT(ITHUMB(1))/10000.
DB(J)=(DP(I)/PARAM(I))+(DP(I+2)/PARAM(I+2))
DY=A1*DP(IP2)/XYLD(ENORM(J))
DY=DY*(XYLD(ENORM(J))-XYLD(ENORM(J)+0.02))*50.
DB(J)=(DB(J)+DY)*BNORMX(J)
CALL TOGGLS(IT)
IF(BITEST(IT,14))WRITE(6,125) J,PARAM(IP2),ANORMX,ENORM(J),
1 BNORMX(J),DB(J),ELE(J)
J=J+1
91 CONTINUE
J=J-1
IF(OUTPT) WRITE(6,109) (I,PARAM(I),DP(I),I=1,NPARAM)
109 FORMAT(1X,2HP(,I2,2H)=,F12.3,5X,6HDELTA=,F12.3)
125 FORMAT(1X,I4,2X,F8.3,F11.6,3X,F8.3,3X,E11.5,2X,E11.5,3X,1A5)
RETURN
END

```

```

C.....FUNCTION XYLD, TO TRANSFORM RATIO TO RHODIUM TO
1 NANOGRAMS/CM**2
FUNCTION XYLD(E)
REAL P(8)
DATA P(1),P(2),P(3),P(4),P(5),P(6),P(7)/
1 1.2347,30.8516,-306.138,1980.348,-6477.202,
2 10661.891,-7126.853/
1 AMT=0.
DO 2 I=1,7
AMT=AMT+P(I)/(E)**(I-1)
2 CONTINUE
XYLD=10.***AMT
RETURN
END

```

```

SUBROUTINE LABEL(J,JPEAK)
REAL ENORM(30), ELE(30), K, MN, CRMN, NI, NIBET, MO, BNORMX(30), DB(30)
INTEGER CH0(30), LIST(520), DATAD(4096)
LOGICAL CHECK
COMMON/FINOUT/ELE,A1,A2
COMMON DATAD
EQUIVALENCE (CH0(1),DATAD(4021)),(LIST(1),DATAD(3073)),
1 (ENORM(1),DATAD(3800)),(BNORMX(1),DATAD(3861)),
2 (DB(1),DATAD(3921))
DATA CL,AR,SC,CA, TI/2HCL,2HAR,1HK,2HSC,2HCA,2HTI/,
1 V,CR,MN,FE,CO,NI/1HV,2HCR,2HMN,0HFE,2HCO,2HNI/,
2 CU,ZN,TLL,PBL,SE,BR/2HCU,2HZN,3HTLL,3HPRL,2HSE,2HBF/,
3 THL,RB,SR,ZR,MO/3HTHL,2HRR,2HSR,2HZR,2HMO/,
4 CASC,SCTI,TIV,VCR,CRMN/5HCA-SC,5HSC-TI,4HTI-V,
5 4HV-CP,5HCR-MN/,
6 FEBET,NIBET,CUBET/5HFERET,5HNIBET,5HCURET/,
7 ZNBET,BRBET,UNKNW/5HZNRET,5HBRRET,5H----/
NELE=(LIST(6)/3)+J
DO 2 I=J,NELE

```

```

E=ENORM(I)
ELE(I)=UNKNW
IF(ABS(E-2.62).LE.0.06) ELE(I)=CL
IF(ABS(E-2.95).LE.0.06) ELE(I)=AR
IF(ABS(E-3.31).LE.0.06) ELE(I)=K
IF(ABS(E-3.69).LE.0.06) ELE(I)=CA
IF(ABS(E-4.09).LE.0.06) CALL SORT(CHECK,3.69,JPEAK)
IF(.NOT. CHECK.AND.(ABS(E-4.09).LT.0.06)) ELE(I)=SC
IF(CHECK) ELE(I)=CASC
CHECK=.FALSE.
IF(ABS(E-4.51).LE.0.06) CALL SORT(CHECK,4.09,JPEAK)
IF(.NOT.CHECK.AND.(ABS(E-4.51).LT.0.06)) ELE(I)=TI
IF(CHECK) ELE(I)=SCTI
CHECK=.FALSE.
IF(ABS(E-4.95).LE.0.06) CALL SORT(CHECK,4.51,JPEAK)
IF(.NOT.CHECK.AND.(ABS(E-4.95).LE.0.06)) ELE(I)=V
IF(CHECK) ELE(I)=TIV
CHECK=.FALSE.
IF(ABS(E-5.41).LE.0.06) CALL SORT(CHECK,4.95,JPEAK)
IF(.NOT.CHECK.AND.(ABS(E-5.41).LE.0.06)) ELE(I)=CR
IF(CHECK) ELE(I)=VCR
CHECK=.FALSE.
IF(ABS(E-5.90).LE.0.06) CALL SORT(CHECK,5.41,JPEAK)
IF(.NOT.CHECK.AND.(ABS(E-5.90).LE.0.06)) ELE(I)=MN
IF(CHECK) ELE(I)=CRMN
CHECK=.FALSE.
28 IF(ABS(E-6.40).LE.0.06) ELE(I)=FE
IF(ABS(E-6.93).LE.0.06) ELE(I)=CO
IF(ABS(E-7.06).LE.0.06) ELE(I)=FERET
IF(ABS(E-7.48).LE.0.06) ELE(I)=NI
IF(ABS(E-8.05).LE.0.06) ELE(I)=CU
IF(ABS(E-8.26).LE.0.06) CALL SORT(CHECK,7.48,JPEAK)
IF(CHECK) ELE(I)=NIBET
CHECK=.FALSE.
IF(ABS(E-8.63).LE.0.06) ELE(I)=ZN
IF(ABS(E-8.90).LE.0.06) CALL SORT(CHECK,8.05,JPEAK)
IF(CHECK) ELE(I)=CUBET
CHECK=.FALSE.
IF(ABS(E-9.57).LE.0.06) CALL SORT(CHECK,8.63,JPEAK)
IF(CHECK) ELE(I)=ZNBET
CHECK=.FALSE.
IF(ABS(E-10.25).LE.0.06) CALL SORT(CHECK,12.21,JPEAK)
IF(CHECK) ELE(I)=TLL
CHECK=.FALSE.
IF(ABS(E-10.55).LE.0.06) CALL SORT(CHECK,12.61,JPEAK)
IF(CHECK) ELE(I)=PBL
CHECK=.FALSE.
34 IF(ABS(E-11.22).LE.0.06) ELE(I)=SE
IF(ABS(E-11.89).LE.0.06) ELE(I)=BR
IF(ABS(E-12.21).LE.0.06) CALL SORT(CHECK,10.27,JPEAK)
IF(CHECK) ELE(I)=TLL
CHECK=.FALSE.
IF(ABS(E-12.61).LE.0.06) CALL SORT(CHECK,10.55,JPEAK)
IF(CHECK) ELE(I)=PBL
CHECK=.FALSE.
IF(ABS(E-12.90).LE.0.10) CALL SORT(CHECK,16.20,JPEAK)
IF(CHECK) ELE(I)=THL
CHECK=.FALSE.
IF(ABS(E-13.29).LE.0.06) CALL SORT(CHECK,11.92,JPEAK)

```

```

IF(CHECK) ELE(I)=BRRET
CHECK=.FALSE.
IF(ABS(E-13.39).LE.0.06) ELE(I)=RB
IF(ABS(E-14.16).LE.0.06) ELE(I)=SR
IF(ABS(E-15.77).LE.0.06) ELE(I)=ZR
IF(ABS(E-16.20).LE.0.10) CALL SORT(CHECK,12.90,JPEAK)
IF(CHECK) ELE(I)=THL
CHECK=.FALSE.
42 IF(ABS(E-17.47).LE.0.26) ELE(I)=MO
2 CONTINUE
RETURN
END

```

```

SUBROUTINE SORT(CHECK,X,JPEAK)
INTEGER CH0(30),DATAD(4096)
LOGICAL CHECK
REAL ELE(30),BNORMX(30),DB(30)
COMMON/FINOUT/ELE,A1,A2
COMMON/DATAD
EQUIVALENCE (CH0(1),DATAD(4021)),(BNORMX(1),DATAD(3861)),
1 (DB(1),DATAD(3921))
CHECK=.FALSE.
DO 2 I=1,JPEAK
E=A1*FLOAT(CH0(I))+A2
IF(ABS(E-X).LE.0.10) GO TO 5
GO TO 2
5 I=JPEAK
CHECK=.TRUE.
2 CONTINUE
RETURN
END

```

```

SUBROUTINE SCALI(NX,NY,NF)
C.....SUBROUTINE TO CHANGE SCALES ON DISPLAY,
C.....FOR USE WITH IPICK
INTEGER GATE(5,30),DATA(1024),CH0(30),DATAD(4096),CHMAX
COMMON/STUFF/GATE
COMMON/DATAD
EQUIVALENCE (DATA(1),DATAD(1025)),(CH0(1),DATAD(4021))
I2=IROTOR(2)
NF=LIMIT(1,256*I2-255,1024)
I3=IROTOP(3)
NX=LIMIT(8,I3+7,12)
I4=IROTOR(4)
NY=LIMIT(8,I4+6,18)
DO 10 I=1,30
LI=CH0(I)
IF(LI.EQ.0) LI=1
GATE(1,I)=LIMIT(0,(CH0(I)-NF)*8/2**I3,1023)
GATE(2,I)=DATA(LI)/(2**I4)*16
10 CONTINUE
RETURN
END

```

```

SUBROUTINE REED(ICTYPE,FNAME,NEXT,EXIST)
C.....SUBROUTINE TO READ DATA FROM TAPE FOR USE WITH IPK2
C.....AND SIMUL2
      INTEGER DATA(1024),DATAD(4096),BKGND(1024)
      REAL FNAME(2)
      LOGICAL EXIST,BITEST
      COMMON DATAD
      EQUIVALENCE (DATA(1),DATAD(1025)),(BKGND(1),DATAD(2049))
      CALL TOGGLS(IT)
      IF(BITEST(IT,15)) GO TO 3
1     CALL INPUT1(FNAME(1),5HNAME,6)
      CALL INPUT1(NEXT,4HNEXT,1)
3     CALL BCDEXT(FNAME,NEXT)
      CALL FSTAT(1,FNAME,EXIST)
      IF(EXIST) GO TO 2
      WRITE(6,100) FNAME
      IF(BITEST(IT,15))RETURN
      GO TO 1
2     IF(ICTYPE.EQ.1)CALL ZERO(DATA)
      IF(ICTYPE.EQ.1) WRITE(6,101) FNAME
101   FORMAT(1HI,12H PUN NUMBER ,2A5)
      IF(ICTYPE.EQ.2) CALL ZERO(BKGND)
      CALL SEEK(1,FNAME)
      IF(ICTYPE.EQ.1)READ(1) TS,IRT,ILT,ND,(DATA(I),I=1,1024)
      IF(ICTYPE.EQ.2) READ(1) TS,IRT,ILT,ND,(BKGND(I),I=1,1024)
      RETURN
100   FORMAT(1X,6H0FILE ,2A5,12HNOT FOUND.)
      END

```

```

SUBROUTINE MAXER(DATA,CHMAX)
      INTEGER DATA(1024),CHMAX
      CHMAX=450
      DO 400 I=450,1024
      IF(DATA(I).GT.DATA(CHMAX))CHMAX=I
400   CONTINUE
      RETURN
      END

```

```

SUBROUTINE SET1(CHMAX, IDAT, POINT)
C.....ROUTINE TO SET UP CALL TO FITALL TO FIND HEIGHT OF
C       NORMALIZATION PEAK
      INTEGER CHMAX, IDAT(1024), DATAD(4096), LIST(520)
      REAL POINT(100), PARAM(15)
      COMMON DATAD
      EQUIVALENCE (LIST(1),DATAD(3073)),(PARAM(1),DATAD(4056))
      IMAX=CHMAX+20
      IMIN=CHMAX-20
      BG2=IDAT(IMAX)
      BG1=IDAT(IMIN)
      DO 200 I=1,41
      ICO=IMIN+I-1
      POINT(I)=IDAT(ICO)
200   CONTINUE
      PARAM(2)=(BG1)
      PARAM(1)=(BG2-BG1)/40.
      PARAM(5)=6./1.665

```

```

PARAM(3)=FLOAT(IDAT(CHMAX))-PARAM(2)-PARAM(1)*20.
PARAM(4)=FLOAT(CHMAX-IMIN+1)
LIST(1)=41
LIST(2)=5
LIST(3)=15
LIST(4)=0
LIST(5)=5
LIST(6)=5
RETURN
END

SUBROUTINE OUTER(PARAM, LIST, ANORM, ARNORM, CHMAX)
INTEGER CHMAX, LIST(520)
REAL PARAM(15)
PARAM(4)=PARAM(4)+FLOAT(CHMAX-21)
IF(LIST(4).NE.0) WRITE(6,111) LIST(4)
111 FORMAT(1X,1IHEX,TYPE ,I2)
ANORM=PARAM(3)*PARAM(5)*1.7725
ABNORM=40.*PARAM(2)+800.*PARAM(1)
WRITE(6,123) PARAM(4),ANORM
WRITE(6,124) ABNORM
FORMAT(1X,15HBACKGROUND HAS ,F10.3,8H COUNTS.)
FORMAT(1X,29HNORMALIZATION PEAK IN CHANNEL
1L,F8.2,4H HAS,F10.2,8H COUNTS.)
124 RETURN
123 END

SUBROUTINE SUBTR(ANORM, ARNORM, CHMAX, OUTPT, GAM)
INTEGER DATA(1024),CHMAX,DATAD(4096),BKGND(1024)
REAL PARAM(15)
LOGICAL OUTPT,BITEST
COMMON DATAD
EQUIVALENCE (DATA(1),DATAD(1025)),(BKGND(1),DATAD(2049)),
1 (PARAM(1),DATAD(4056))
PARAM(4)=PARAM(4)+FLOAT(CHMAX-21)
CNORM=PARAM(3)*PARAM(5)*1.7725
CBNORM=40.*PARAM(2)+800.*PARAM(1)
WRITE(6,123) PARAM(4),CNORM
WRITE(6,124) CBNORM
FORMAT(1X,15HBACKGROUND HAS ,F10.3,8H COUNTS.)
RAT=ABNORM/CBNORM
DO 5 I=3,1023
IS=I-2
DATA(I)=FLOAT(DATA(I))-RAT*FLOAT(BKGND(I-1)+BKGND(I)+
1 BKGND(I+1))/3.
BKGND(IS)=FLOAT(DATA(I))+2.*RAT*FLOAT(BKGND(I-1)+
1 BKGND(I)+BKGND(I+1))/3.
IF(DATA(I).LT.0) DATA(I)=0
5 CONTINUE
BKGND(1023)=0
BKGND(1024)=0
CALL TOGLLS(IT)
IF(BITEST(IT,15)) RETURN
WRITE(6,105)
CALL INPUTI(OUTPT,5HOUTPT,4)
CALL INPUTI(GAM,5HGAMMA,2)

```

```

123  FORMAT(//,1X,29HNORMALIZATION PEAK IN CHANNEL
1L,F8.2,4H HAS,F10.2,8H COUNTS.)
105  FORMAT(1X,25HDO YOU WISH EXTRA OUTPUT?)
      RETURN
      END

      SUBROUTINE MULT(IMAX,IMIN,J)
      INTEGER CH0(30),DATAD(4096)
      LOGICAL DOUBLE,UNDOUN,TRIPLE,UNTRIP,QUAD,UNQUAD,QUIN,UNQUIN
      COMMON/MUP/ DOUBLE,UNDOUN,TRIPLE,UNTRIP,QUAD,UNQUAD,QUIN,
      1    UNQUIN,JPEAK,
      1    GAM, IDUB, ITRIP, IQUAD, IQUIN
      COMMON DATAD
      EQUIVALENCE (CH0(1),DATAD(4021))
      DOUBLE=.FALSE.
      TRIPLE=.FALSE.
      QUAD=.FALSE.
      QUIN=.FALSE.
      I0=CH0(J)
      IMAX=CH0(J)+IFIX(GAM)
      IMIN=CH0(J)-IFIX(GAM)
      IF(UNDOUN) GO TO 82
      IF(J.EQ.JPEAK) GO TO 82
      IF(CH0(J+1).LT.(I0+IFIX(4.*GAM))) DOUBLE=.TRUE.
      IDUB=CH0(J+1)
      IF(UNTRIP) GO TO 82
      IF(J.EQ.JPEAK-1) GO TO 82
      IF(CH0(J+2).LT.(IDUB+IFIX(4.*GAM))) TRIPLE=.TRUE.
      TRIPLE=DOUBLE.AND.TRIPLE
      ITRIP=CH0(J+2)
      IF(UNQUAD) GO TO 82
      IF(J.EQ.JPEAK-2) GO TO 82
      IF(CH0(J+3).LT.(ITRIP+IFIX(4.*GAM))) QUAD=.TRUE.
      QUAD=TRIPLE.AND.QUAD
      IQUAD=CH0(J+3)
      IF(UNQUIN) GO TO 82
      IF(J.EQ.JPEAK-3) GO TO 82
      IF(CH0(J+4).LT.(IQUAD+IFIX(4.*GAM))) QUIN=.TRUE.
      QUIN=QUIN.AND.QUAD
      IQUIN=CH0(J+4)
82   UNDOUN=.FALSE.
      UNTRIP=.FALSE.
      UNQUAD=.FALSE.
      UNQUIN=.FALSE.
      IF(DOUBLE) IMAX=IDUB+IFIX(GAM)
      IF(TRIPLE) IMAX=ITRIP+IFIX(GAM)
      IF(QUAD) IMAX=IQUAD+IFIX(GAM)
      IF(QUIN) IMAX=IQUIN+IFIX(GAM)
      RETURN
      END

      SUBROUTINE SIT2(IMAX,IMIN,J,ANORM)
C.....TO PREPARE FINAL FIT TO DATA
      REAL PARAM(15),POINT(100)
      INTEGER DATA(1024),LIST(520),CH0(30),IR(5),DATAD(4096)
      LOGICAL DOUBLE,UNDOUN,TRIPLE,UNTRIP,QUAD,UNQUAD,QUIN,UNQUIN,
      1    BITEST

```

```

COMMON/MUP/DOUBLE, UNDOUB, TRIPLE, UNTRIP, QUAD, UNQUAD, QUIN,
1      UNQUIN, JPEAK, GAM, IDUB, ITRIP, IQUAD, IQUIN
COMMON DATAD
EQUIVALENCE (DATA(1),DATAD(1025)),(LIST(1),DATAD(3073)),
1      (CH0(1),DATAD(4021)),(POINT(1),DATAD(3593)),
2      (IR(1),DATAD(4051)),(PARAM(1),DATAD(4056))
ITOT=IMAX-IMIN+1
DO 84 I=1,ITOT
IDOWN=IMIN+I-1
POINT(I)=FLOAT(DATA(IDOWN))
84  CONTINUE
I0=CH0(J)
PARAM(14)=FLOAT(IQUIN-IMIN+1)
PARAM(15)=GAM/1.665
PARAM(13)=FLOAT(DATA(IQUIN))
PARAM(11)=FLOAT(IQUAD-IMIN+1)
PARAM(12)=GAM/1.665
PARAM(10)=FLOAT(DATA(IQUAD))
85  PARAM(7)=FLOAT(DATA(ITRIP))
PARAM(8)=FLOAT(ITRIP-IMIN+1)
PARAM(9)=GAM/1.665
86  PARAM(4)=FLOAT(DATA(IDUB))
PARAM(5)=FLOAT(IDUB-IMIN+1)
PARAM(6)=GAM/1.665
87  PARAM(1)=FLOAT(DATA(I0))
PARAM(2)=FLOAT(I0-IMIN+1)
PARAM(3)=GAM/1.665
LIST(1)=ITOT
LIST(2)=3
LIST(3)=15
LIST(4)=0
LIST(5)=12
LIST(6)=3
IF(.NOT.DOUBLE) GO TO 88
LIST(2)=6
LIST(6)=6
IF(.NOT.TRIPLE) GO TO 88
LIST(2)=9
LIST(6)=9
IF(.NOT.QUAD) GO TO 88
LIST(2)=12
LIST(6)=12
IF(.NOT.QUIN) GO TO 88
LIST(2)=15
LIST(6)=15
LIST(5)=15
88  CONTINUE
NPARAM=LIST(6)
CALL ZERO(IR)
ISUM=0
K=0
DO 95 I=1,NPARAM,3
K=K+1
IF(PARAM(I).LT.0.0000300*ANORM) IR(K)=1
ISUM=ISUM+IR(K)
95  CONTINUE
LIST(2)=LIST(2)-3*ISUM
LIST(5)=LIST(5)-ISUM*3
LIST(6)=LIST(6)-3*ISUM

```

```

11=1
DO 97 I=1,K
IF(IRC(I).NE.0) GO TO 97
96 PARAM(I1)=PARAM(3*I-2)
PARAM(I1+1)=PARAM(3*I-1)
PARAM(I1+2)=PARAM(3*I)
98 I1=I1+3
CONTINUE
97 CALL TOGLLS(IT)
IF(DOUBLE.AND.(.NOT.TRIPLE).AND.BITEST(IT,14)) WRITE(6,115)
IF(TRIPLE.AND.(.NOT.QUAD).AND.BITEST(IT,14)) WRITE(6,116)
IF(QUAD.AND.(.NOT.QUIN).AND.BITEST(IT,14)) WRITE(6,118)
IF(QUIN.AND.BITEST(IT,14)) WRITE(6,119)
RETURN
115 FORMAT(1X,8H DOUBLET)
116 FORMAT(1X,8H TRIPLET)
118 FORMAT(1X,11H QUADRUPLET)
119 FORMAT(1X,11H QUINTUPLET)
END

```

```

SUBROUTINE IDENT(ELE,ENORM,BNORMX,DR,JPEAK,FNAME,NEXT)
C.....SUBROUTINE TO MAKE FINAL IDENTIFICATION OF ELEMENTS,
C.....AND MAKE CORRECTIONS FOR BETA PEAKS.
REAL ELE(30),ENORM(30),BNORMX(30),DR(30),FNAME(2),
1 ESI(24),ES2(24),ATOR(24),NAM(24),EFF(4)
LOGICAL BITEST,BETA
DATA ESI(1),ES2(1),ATOR(1),NAM(1)/2.62,2.81,0.,2HCL/,
1 ESI(2),ES2(2),ATOR(2),NAM(2)/2.96,3.19,0.,2HAR/,
2 ESI(3),ES2(3),ATOR(3),NAM(3)/3.313,3.589,0.114,1HK/,
3 ESI(4),ES2(4),ATOR(4),NAM(4)/3.691,4.012,0.0909,2HCA/,
4 ESI(5),ES2(5),ATOR(5),NAM(5)/4.090,4.460,0.105,2HSC/,
5 ESI(6),ES2(6),ATOR(6),NAM(6)/4.510,4.931,0.105,2HTI/,
6 ESI(7),ES2(7),ATOR(7),NAM(7)/4.952,5.427,0.105,1HV/
DATA ESI(8),ES2(8),ATOR(8),NAM(8)/5.414,5.946,0.114,2HCP/,
1 ESI(9),ES2(9),ATOR(9),NAM(9)/5.898,6.490,0.108,2HMN/,
2 ESI(10),ES2(10),ATOR(10),NAM(10)/6.403,7.057,0.106,2HFE/,
3 ESI(11),ES2(11),ATOR(11),NAM(11)/6.930,7.649,0.118,2HCO/,
4 ESI(12),ES2(12),ATOR(12),NAM(12)/7.477,8.264,0.105,2HNI/,
5 ESI(13),ES2(13),ATOR(13),NAM(13)/8.047,8.904,0.113,2HCU/
DATA ESI(14),ES2(14),ATOR(14),NAM(14)/8.638,9.571,0.117,2HZN/,
1 ESI(15),ES2(15),ATOR(15),NAM(15)/10.543,11.725,0.083,2HAS/,
2 ESI(16),ES2(16),ATOR(16),NAM(16)/11.221,12.495,0.13,2HSE/,
3 ESI(17),ES2(17),ATOR(17),NAM(17)/11.923,13.290,0.136,2HBR/,
4 ESI(18),ES2(18),ATOR(18),NAM(18)/14.164,15.834,0.151,2HSR/,
5 ESI(19),ES2(19),ATOR(19),NAM(19)/15.774,17.666,0.172,2HZR/
DATA ESI(20),ES2(20),ATOR(20),NAM(20)/17.478,19.607,0.18,2HMO/,
1 ESI(21),ES2(21),ATOR(21),NAM(21)/8.396,9.670,0.775,1HW/,
2 ESI(22),ES2(22),ATOR(22),NAM(22)/10.266,12.210,0.672,2HTL/,
3 ESI(23),ES2(23),ATOR(23),NAM(23)/10.549,12.611,0.645,2HPB/,
4 ESI(24),ES2(24),ATOR(24),NAM(24)/12.966,16.200,0.553,2HTH/
DATA EPS,UNKNW/0.060,2H--/
DATA EFF(1),EFF(2),EFF(3),EFF(4)/1.775,1.66,1.66,2.00/
MEXT=NEXT-1
CALL BCDEXT(FNAME,MEXT)
DO 4 I=3,JPEAK
ELE(I)=UNKNW
DO 2 J=1,24
IF(ABS(ENORM(I)-ESI(J)).LT.EPS) GO TO 1

```

```

1      GO TO 2
BETA=.FALSE.
DO 3 K=I,JPEAK
IF(ABS(ENORM(K)-ES2(J)).GT.0.090) GO TO 3
IF(BNORMX(K).GT.0.3*BNORMX(I)) BETA=.TRUE.
BNORMX(K)=BNORMX(K)-ATOR(J)*BNORMX(I)
IF(BNORMX(K).LT.0.) BNORMX(K)=0.
3      CONTINUE
IF(J.LT.21) ELE(I)=NAM(J)
IF(J.GE.21 .AND. BETA) GO TO 6
GO TO 2
6      ELE(I)=NAM(J)
BNORMX(I)=BNORMX(I)*EFF(J-20)
DR(I)=DB(I)*EFF(H-20)
2      CONTINUE
4      CONTINUE
CALL TOGGLS(IT)
IF(.NOT.BITEST(IT,13)) GO TO 5
CALL ENTER(3,FNAME)
100   WRITE(3,100) (ELE(I),BNORMX(I),DR(I),I=3,JPEAK)
FORMAT(1X,1A5,2E11.5)
CALL CLOSE(3)
5      CALL TOGGLS(IT)
IF(.NOT.BITEST(IT,12)) RETURN
WRITE(6,103) FNAME
C..... WRITE(6,104) G
C..... FORMAT(1X,2HG=,F6.4)
WRITE(6,101)
WRITE(6,102) (ELE(I),BNORMX(I),DR(I),I=3,JPEAK)
RETURN
101   FORMAT(1X,7HELEMENT,2X,13HCONCENTRATION,9H      ERROR)
102   FORMAT(4X,1A5,E12.6,E12.6)
103   FORMAT(1H1,5X,2A5)
END

```

```

SUBROUTINE ERRO(J,IMIN,DP)
REAL ELE(30),BNORMX(30),DP(15),PARAM(15)
INTEGER LIST(520),DATAD(4096),IR(5),CH0(30)
LOGICAL DOUBLE,UNDOUB,TRIPLE,UNTRIP,QUAD,UNQUAD,QUIN,UNQUIN,
I      BITEST
COMMON /MUP/DOUBLE,UNDOUB,TRIPLE,UNTRIP,QUAD,UNQUAD,QUIN,
I      UNQUIN,JPEAK,GAM,IDUB,ITRIP,IQUAD
COMMON DATAD
COMMON/FINOUT/ELE,A1,A2
EQUIVALENCE (LIST(1),DATAD(3073)),(BNORMX(1),DATAD(3861)),
1      (DB(1),DATAD(3921)),(IR(1),DATAD(4051)),(DATAD(4021),
2      CH0(1)),(PARAM(1),DATAD(4056))
DATA RNONE/4HNONE/
CALL TOGGLS(IT)
IF(BITEST(IT,14))WRITE(6,111) LIST(4)
111   FORMAT(1H ERROR,TYPE,I2)
IF(.NOT.DOUBLE) GO TO 89
IF(.NOT.TRIPLE) GO TO 90
IF(.NOT.QUAD) GO TO 91
IF(.NOT.QUIN) GO TO 92
UNQUIN=.TRUE.
RETURN

```

```

92      UNQUAD=.TRUE.
91      RETURN
90      UNTRIP=.TRUE.
90      RETURN
89      UNDOUB=.TRUE.
89      RETURN
89      PARAM(1)=0.
89      PARAM(3)=0.
89      PARAM(2)=FLOAT(CH0(J)-IMIN+1)
89      DP(1)=0.
89      DP(2)=0.
89      DP(3)=0.
89      RETURN
89      END

SUBROUTINE ENERG(IMIN,A1,A2,J,DP)
REAL PARAM(15),ENORM(30),DP(15)
INTEGER DATAD(4096),LIST(520),IR(5),CH0(30)
COMMON DATAD
EQUIVALENCE (DATAD(3800),ENORM(1)),(LIST(1),DATAD(3073)),
1      (CH0(1),DATAD(4021)),(IR(1),DATAD(4051)),
2      (PARAM(1),DATAD(4056))
K=J
ISUM=0
DO 2 I=1,5
IF(IR(I).EQ.2) GO TO 2
ISUM=ISUM+IR(I)
2 CONTINUE
K1=0
NPPEAK=LIST(6)/3
NPEAK=NPPEAK+ISUM
DO 3 I=1,NPEAK
MPEAK=NPEAK-I+1
M3PEAK=3*MPEAK
IF(IR(MPEAK).EQ.0) GO TO 4
JMPEAK=J+MPEAK-1
PARAM(M3PEAK)=0.
DP(M3PEAK)=0.
PAPAM(M3PEAK-1)=FLOAT(CH0(JMPEAK))-FLOAT(IMIN-1)
DP(M3PEAK-1)=0.
PARAM(M3PEAK-2)=0.
DP(M3PEAK-2)=0.
GO TO 3
4 N3PEAK=3*(NPPEAK-K1)
PARAM(M3PEAK)=PARAM(N3PEAK)
DP(M3PEAK)=DP(N3PEAK)
PAPAM(M3PEAK-1)=PARAM(N3PEAK-1)
DP(M3PEAK-1)=DP(N3PEAK-1)
PAPAM(M3PEAK-2)=PARAM(N3PEAK-2)
DP(M3PEAK-2)=DP(N3PEAK-2)
K1=K1+1
3 CONTINUE
LIST(6)=LIST(6)+3*ISUM
NPARAM=LIST(6)
DO 5 I=1,NPARAM,3
IP2=I+1
PARAM(IP2)=PARAM(IP2)+FLOAT(IMIN-1)
ENORM(K)=A1*PARAM(IP2)+A2

```

```

      K=K+1
5    CONTINUE
      RETURN
      END

      SUBROUTINE LOCUM
C.....TO LOCATE PEAKS IN X-RAY FLUORESCENT ENERGY SPECTRA..
C.....TO BE USED WITH IPK3, AS PART OF AN EXECUTE FILE..SAVLOC
      INTEGER DATA(1024),GATE(5,30),CH0(30),DATAD(4096),CHMAX,
     1   CH1,CH2,CH3,CH4,CH5,BKGND(1024),GATI(140)
      REAL C(36),B(36),PHI2,PHI
      LOGICAL DOUBLE,UNDOUB,TRIPLE,UNTRIP,QUAD,UNQUAD,
     1   QUIN,UNQUIN
      COMMON/STUFF/GATE
      COMMON/MUP/DOUBLE,UNDOUB,TRIPLE,UNTRIP,QUAD,UNQUAD,
     1   QUIN,UNQUIN,JPEAK,GAM,IDUB,ITRIP,IQUAD,IQUIN
      COMMON DATAD
      COMMON/DIFF/C,B,M
      EQUIVALENCE (DATA(1),DATAD(1025)),(CH0(1),DATAD(4021)),
     1   (BKGND(1),DATAD(2049)),(GATI(1),GATE(1,3))
      CALL ZERO(GATI)
      NHMAX=470
      M=IFIX((0.6*GAM-1.)/2.)
      IF(M.EQ.0) M=1
C.....COMPUTE COEFFICIENT C(J)
      CALL ZERO(C)
      CALL ZERO(B)
      C(1)=-0.2
      C(2)=0.1
      C(3)=0.
      DO 31 L=1,5
      DO 30 J=1,36
      B(J)=C(J)
      DO 30 K=1,M
      K1=J+K
      K2=J-K
      IF(J-K.LE.0) K2=K-J+2
      B(J)=B(J)+C(K1)+C(K2)
      IF(K1.GT.36) B(J)=0
30    CONTINUE
      DO 31 J=1,36
      C(J)=B(J)
31    CONTINUE
C.....COMPUTE GENERALIZED SECOND DIFFERENCE, S
C.....COMPUTE STANDARD DEVIATIONS IN S
      PHI=C(1)**2
      M1=5*M+2
      DO 50 J=2,M1
      PHI=PHI+2.*C(J)**2
50    CONTINUE
      PHI2=SQRT(PHI)
C.....COMPUTE CRITERIA FOR PEAK LOCATION
C.....NI=EXPECTED NUMBER OF CHANNELS IN NEGATIVE PEAK
      NI=IFIX(1.22*(GAM+0.5))
      NI MAX=NI+4
      NI MIN=NI-2
C.....BEGIN TO SEARCH FOR PEAKS
      CH5=25

```

```

JPEAK=2
58 JPEAK=JPEAK+1
59 ICHNL=CH5
60 IF(ICHLN.EQ.NHMAX) GO TO 70
ICHLN=ICHLN+1
IF(S(ICHLN).GE.-1.6*PHI2*SQRT(FLOAT(BKGND(ICHLN-2)+1))) GO TO 60
GO TO 62
61 IF(ICHLN.EQ.NHMAX) GO TO 70
ICHLN=ICHLN+1
62 IF(S(ICHLN+1).LT.S(ICHLN).OR.S(ICHLN-1).LT.S(ICHLN)) GO TO 61
IF(S(ICHLN).LT.-1.6*PHI2*SQRT(FLOAT(BKGND(ICHLN-2)+1))) GO TO 62
CH5=ICHLN
GO TO 59
625 CH4=ICHLN
GO TO 64
63 IF(ICHLN.EQ.NHMAX) GO TO 70
ICHLN=ICHLN+1
64 IF(S(ICHLN).LE.0.8*S(CH4)) GO TO 63
CH5=ICHLN-1
ICHLN=CH4
65 IF(ICHLN.EQ.1) GO TO 59
ICHLN=ICHLN-1
IF(S(ICHLN).LT.0.) GO TO 65
CH3=ICHLN+1
GO TO 67
66 IF(ICHLN.EQ.1) GO TO 59
ICHLN=ICHLN-1
67 IF(S(ICHLN).LT.PHI2*SQRT(FLOAT(BKGND(ICHLN-2)+1))) GO TO 66
CH2=ICHLN
GO TO 69
68 IF(ICHLN.EQ.1) GO TO 59
ICHLN=ICHLN-1
69 IF(S(ICHLN).GT.PHI2*SQRT(FLOAT(BKGND(ICHLN-2)+1))) GO TO 68
CH1=ICHLN+1
C.....COMPUTE EXPECTED WIDTHS
N2 MIN=IFIX((PHI2*SQRT(FLOAT(DATA(CH4)+1))/ABS(S(CH4)))
1 *0.5*FLOAT(NI MIN)+0.5)
N2 MAX=IFIX((PHI2*SQRT(FLOAT(DATA(CH4)+1))/ABS(S(CH4)))
1 *0.5*FLOAT(NI MAX)+0.5)
N3 MIN=IFIX(FLOAT(NI MIN)*(1.-2.*(PHI2*SQRT(FLOAT(DATA(CH4)+1))/
1 ABS(S(CH4)))))+0.5)
N3 MAX=IFIX(FLOAT(NI MAX)*(1.-2.*(PHI2*SQRT(FLOAT(DATA(CH4)+1))/
1 ABS(S(CH4)))))+0.5)
C.....IF(CH5-CH3+1.LT.NI MIN) GO TO 59
C.....IF(CH5-CH3+1.GT.NI MAX) GO TO 59
C.....IF(N2 MAX.EQ.0) GO TO 72
C.....IF(CH3-CH2-1.GT.N2 MAX) GO TO 59
GO TO 73
72 IF(CH3-CH2-1.GT.1) CONTINUE
73 IF(CH2-CH1+1.LT.N3 MIN) CONTINUE
CH0(JPEAK)=CH4
GATE(1,JPEAK)=CH4
GATE(2,JPEAK)=DATA(CH4)
GATE(3,JPEAK)=0
GATE(4,JPEAK)=128
GATE(5,JPEAK)=-10
CALL SCALI(NX,NY,NF)
GO TO 58
70 JPEAK=JPEAK-1
RETURN

```

REFERENCES

1. Costrell, Louis, ed.: CAMAC: A Modular Instrumentation System for Data Handling. TID-25875, Atomic Energy Commission, 1972.
2. Bercaw, Robert W.; Fessler, Theodore E.; and Arnold, Jeffrey M.: A Programmable Computer Interface for CAMAC. NASA TN D-7148, 1973.
3. Mariscotti, M. A.: A Method for Automatic Identification of Peaks in the Presence of Background and Its Application to Spectrum Analysis. Nuclear Instruments and Methods., vol. 50, no. 2, May 1967, pp. 309-320.
4. CHAIN AND EXECUTE Utility Programs. DEC-15-YWZA-DNZ, Digital Equipment Corp., 1970.

TABLE I. - ELEMENTS RECOGNIZED BY SAVLOC

(a) K series

Element	K_{α} energy, keV
Chlorine	2. 62
Argon	2. 96
Potassium	3. 313
Calcium	3. 691
Scandium	4. 090
Titanium	4. 510
Vanadium	4. 952
Chromium	5. 414
Manganese	5. 898
Iron	6. 403
Cobalt	6. 930
Nickel	7. 477
Copper	8. 047
Zinc	8. 638
Arsenic	10. 543
Selenium	11. 221
Bromine	11. 923
Strontium	14. 164
Zirconium	15. 774
Molybdenum	17. 478

(b) L series

Element	L_{α} energy, keV	L_{β} energy, keV
Tungsten	8. 396	9. 670
Thallium	10. 266	12. 210
Lead	10. 549	12. 611
Thorium	12. 966	16. 200

TABLE II. - PARAMETERS USED TO
DESCRIBE X-RAY EFFICIENCY
FUNCTION F(E)

I	P(I)
1	0. 855
2	54. 874
3	-705. 322
4	5499. 011
5	-22895. 21
6	48343. 26
7	-40357. 17

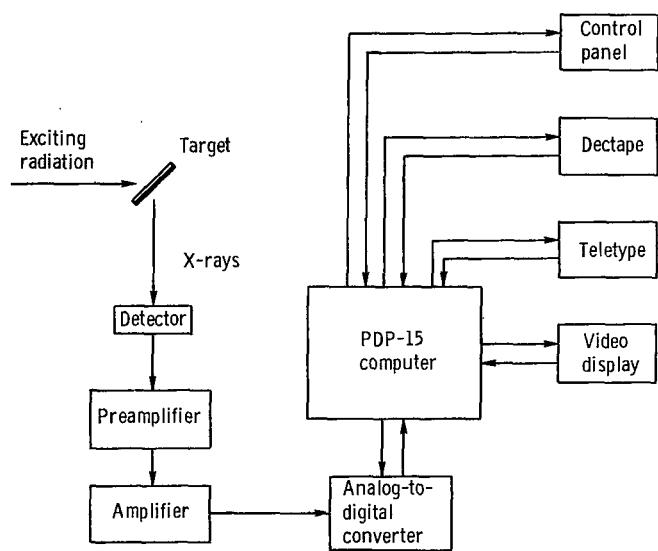


Figure 1. - X-ray fluorescence facility.

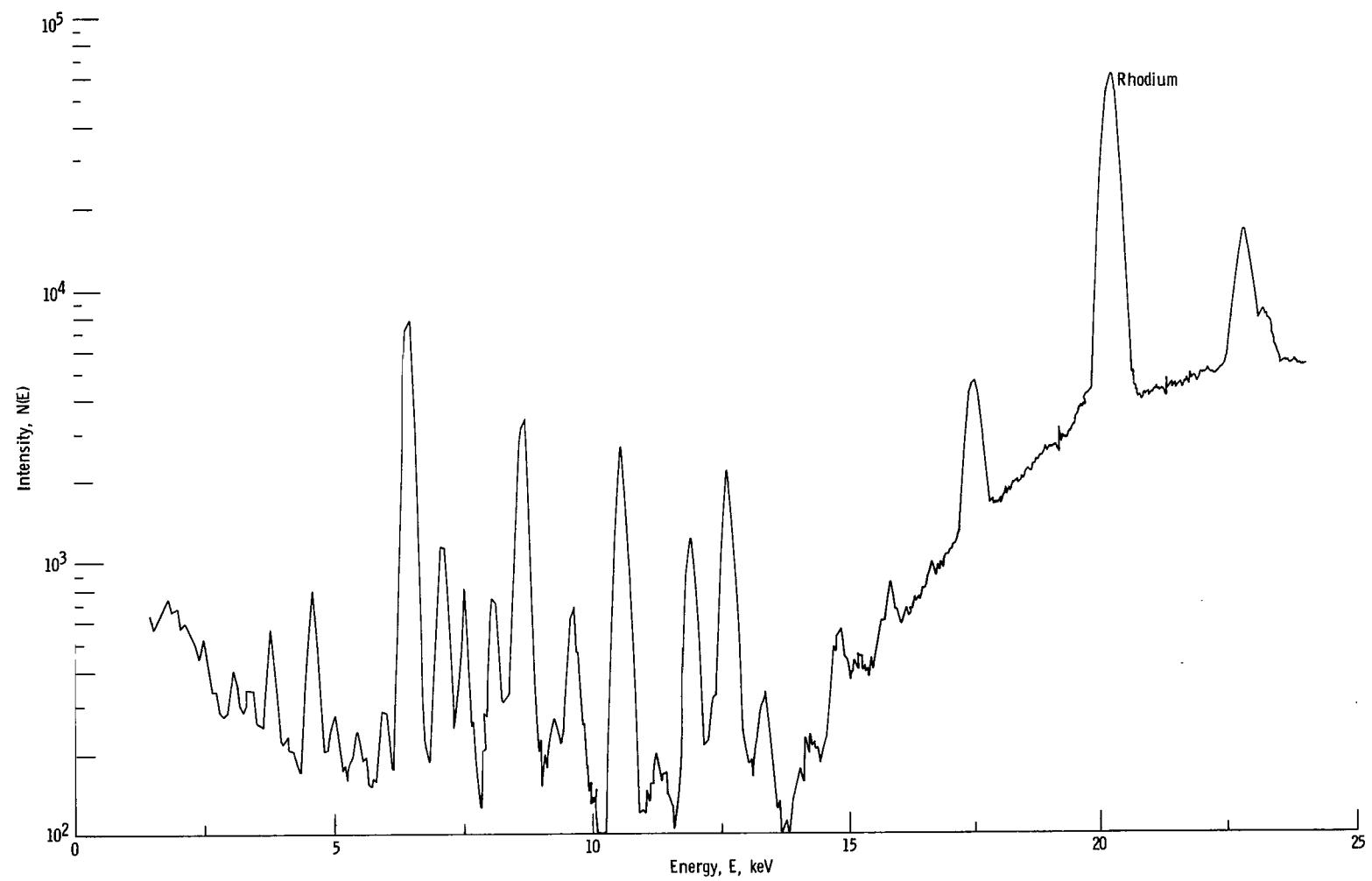


Figure 2. - Typical X-ray spectrum.

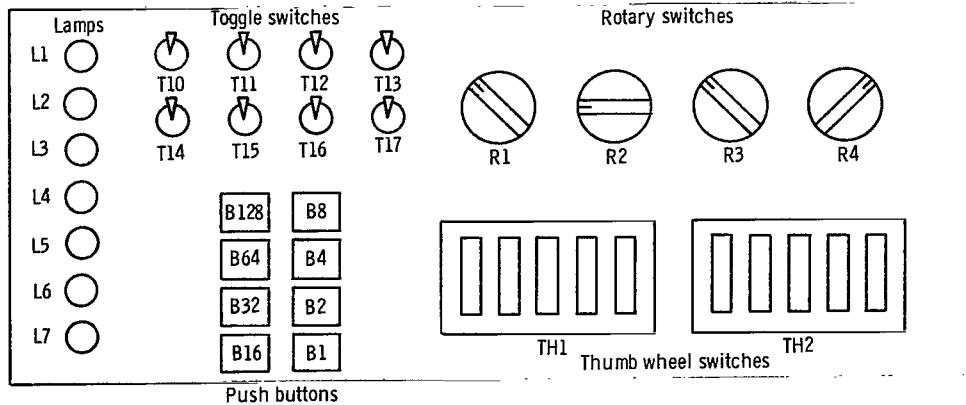


Figure 3. - Functional control panel.

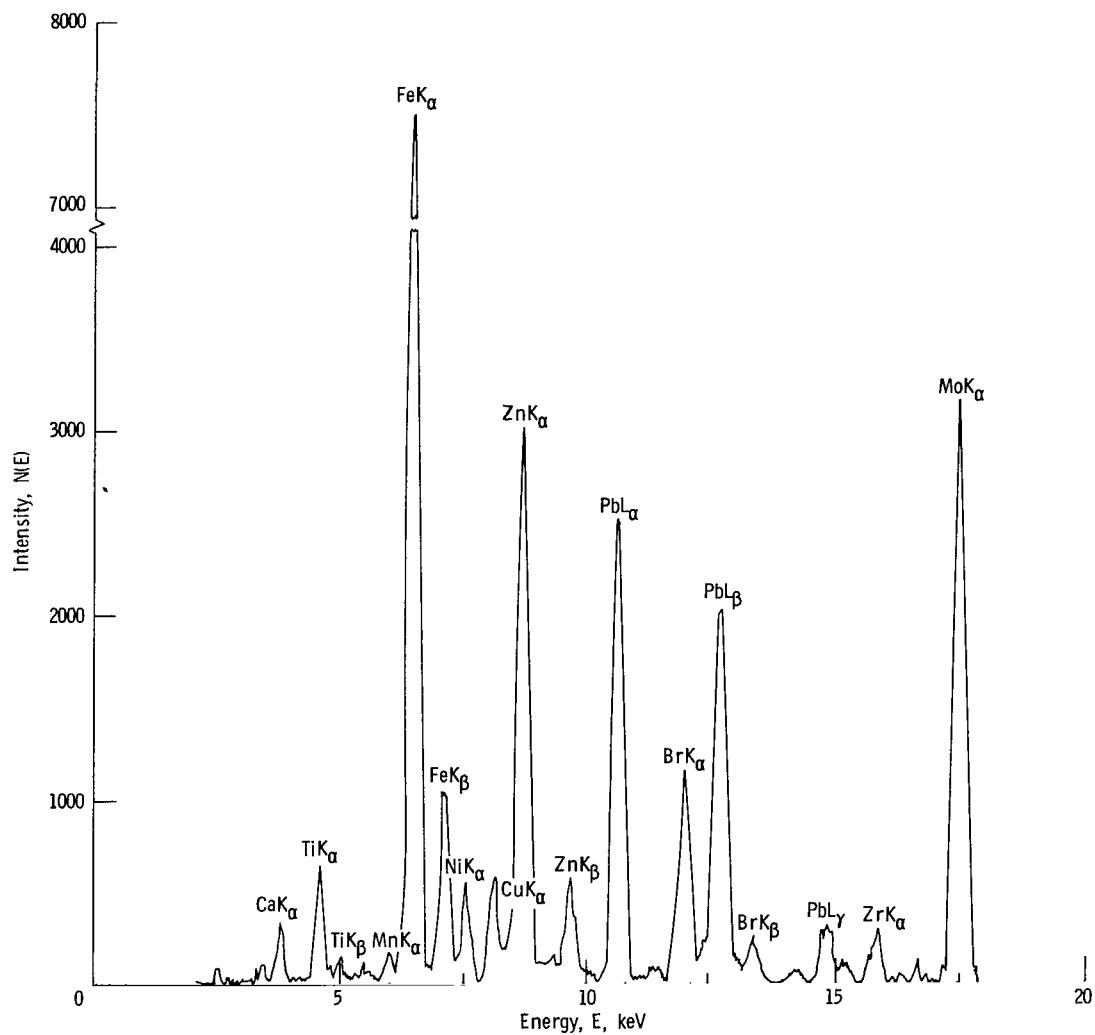


Figure 4. - Typical X-ray spectrum after background subtraction.

RUN NUMBER SEPT 011
 NORMALIZATION PEAK IN CHANNEL .500.93 HAS 526878.78 COUNTS.
 BACKGROUND HAS 152552.379 COUNTS.
 DNAME(A) =SEPT
 NEXT (I) =0

NORMALIZATION PEAK IN CHANNEL .501.48 HAS 648537.67 COUNTS.
 BACKGROUND HAS 56075.779 COUNTS.
 DO YOU WISH EXTRA OUTPUT?
 OUTPT(L) =F
 GAMMA(R) =4.
 ENERGIES FOR CALIBRATION:
 E1 (R) =6.403
 E2 (R) =12.611
 G=1.0000

PEAK	CHANNEL	STRENGTH	ENERGY	CONCENTRATION (UGRAMS/IN**2)	ERROR	ELEM
3	70.035	.000837	3.701	.29701+02	.19374+01	CA
4	91.404	.004320	4.516	.49822+02	.30408+01	TI
DOUBLET						
5	128.004	.000599	5.912	.22555+01	.78361+00	MN
6	140.871	.077698	6.403	.22115+03	.24637+01	FE
QUADRUPLET						
7	157.800	.010517	7.048	.22002+02	.15192+01	FEBET
8	168.506	.001316	7.461	.23297+01	.94050+00	NI
9	183.969	.004440	8.047	.64054+01	.82467+00	CU
10	199.256	.032763	8.630	.39711+02	.13505+01	ZN
11	224.572	.004366	9.595	.41648+01	.61117+00	ZNBET
12	249.423	.029359	10.543	.23081+02	.55729+00	PBL
13	285.057	.012955	11.982	.80868+01	.28089+00	BR
14	303.631	.025692	12.611	.14433+02	.34989+00	PBL
15	321.212	.001912	13.282	.97893+00	.19976+00	BBET
DOUBLET						
16	359.355	.003763	14.737	.16030+01	.17434+00	----
17	368.690	.001194	15.093	.48783+00	.15608+00	----
18	385.632	.003646	15.739	.13832+01	.25890+00	ZR
19	407.321	.000938	16.566	.32513+00	.71051-01	----
20	429.751	.051486	17.422	.16323+02	.29195+00	MO

Figure 5. - Sample of first page of teletype output.

SEPT 011
 ELEMENT CONCENTRATION ERROR
 CA .297007+02 .193737+01
 TI .498224+02 .304084+01
 MN .225552+01 .783605+00
 FE .220910+03 .246371+01
 -- .000000+00 .151919+01
 NI .232971+01 .940501+00
 CU .640538+01 .824675+00
 ZN .397109+02 .135047+01
 -- .000000+00 .611174+00
 PB .341605+02 .824787+00
 BR .808675+01 .280804+00
 -- .000000+00 .349895+00
 -- .000000+00 .199760+00
 -- .162305+01 .174340+00
 -- .487833+00 .156084+00
 ZR .138316+01 .258902+00
 -- .325127+00 .710506-01
 MO .163229+02 .291951+00

Figure 6. - Sample of second page of teletype output.

RUN NUMBER SEPT 011
 NORMALIZATION PEAK IN CHANNEL .500.93 HAS 526878.78 COUNTS.
 BACKGROUND HAS 152552.379 COUNTS.
 DNAME(A) =SEPT
 NEXT (I) =0

NORMALIZATION PEAK IN CHANNEL .501.48 HAS 648537.67 COUNTS.
 BACKGROUND HAS 56075.779 COUNTS.
 DO YOU WISH EXTRA OUTPUT?
 OUTPT(L) =T
 GAMMA(R) =A.
 ENERGIES FOR CALIBRATION:
 E1 (R) =6.403
 E2 (R) =12.611
 G=1.0000

PEAK	CHANNEL	STRENGTH	ENERGY	CONCENTRATION (UGRAMS/IN**2)	ERROR	ELEM
3	70.035	.000837	3.701	.29701+02	.19374+01	CA
P(1)=	213.043		DELTA=	6.546		
P(2)=	70.055		DELTA=	.029		
P(3)=	1.167		DELTA=	.938		
4	91.404	.004320	4.516	.49822+02	.30408+01	TI
P(1)=	509.417		DELTA=	14.783		
P(2)=	91.404		DELTA=	.955		
P(3)=	2.521		DELTA=	.075		
DOUBLET						
5	128.004	.000599	5.912	.22555+01	.78361+00	MN
6	140.871	.077698	6.403	.22115+03	.24637+01	FE
P(1)=	98.433		DELTA=	15.671		
P(2)=	128.004		DELTA=	.235		
P(3)=	1.867		DELTA=	.330		
P(4)=	7328.194		DELTA=	44.861		
P(5)=	140.871		DELTA=	.812		
P(6)=	3.152		DELTA=	.015		
QUADRUPLET						
7	157.800	.010517	7.048	.22002+02	.15192+01	FEBET
8	168.606	.001316	7.461	.23297+01	.94050+00	NI
9	183.969	.004440	8.047	.64054+01	.82467+00	CU
10	199.256	.032763	8.630	.39711+02	.13505+01	ZN
P(1)=	959.873		DELTA=	32.626		
P(2)=	157.800		DELTA=	.079		
P(3)=	3.257		DELTA=	.110		
P(4)=	P 211.486		DELTA=	41.329		
P(5)=	168.606		DELTA=	.280		
P(6)=	1.849		DELTA=	.378		
P(7)=	437.471		DELTA=	28.286		
P(8)=	183.969		DELTA=	.142		
P(9)=	3.017		DELTA=	.188		
P(10)=	2772.324		DELTA=	48.293		
P(11)=	199.256		DELTA=	.042		
P(12)=	3.513		DELTA=	.057		
11	224.572	.004366	9.595	.41648+01	.61117+00	ZNBET
P(1)=	426.654		DELTA=	29.386		
P(2)=	224.572		DELTA=	.157		
P(3)=	3.042		DELTA=	.233		
12	249.423	.029359	10.543	.23081+02	.55729+00	PBL
P(1)=	2379.978		DELTA=	27.044		
P(2)=	249.423		DELTA=	.029		
P(3)=	3.667		DELTA=	.046		
13	285.057	.012955	11.902	.80868+01	.28080+00	BR
P(1)=	1016.039		DELTA=	16.254		
P(2)=	285.057		DELTA=	.042		
P(3)=	3.790		DELTA=	.070		
14	303.631	.025692	12.611	.14433+02	.34989+00	PBL
P(1)=	1891.706		DELTA=	20.430		
P(2)=	303.631		DELTA=	.051		
P(3)=	4.037		DELTA=	.054		
15	321.212	.001912	13.282	.97893+00	.19976+00	BRBET
P(1)=	165.905		DELTA=	14.989		
P(2)=	321.212		DELTA=	.245		
P(3)=	3.428		DELTA=	.385		
DOUBLET						
16	359.355	.003763	14.737	.16030+01	.17434+00	-----
17	368.690	.001194	15.093	.48783+00	.15608+00	-----
P(1)=	294.605		DELTA=	13.175		
P(2)=	359.355		DELTA=	.150		
P(3)=	3.797		DELTA=	.241		
P(4)=	99.216		DELTA=	13.019		
P(5)=	368.690		DELTA=	.429		
P(6)=	3.579		DELTA=	.669		
18	385.632	.003646	15.739	.13832+01	.25890+00	ZR
P(1)=	242.082		DELTA=	17.146		
P(2)=	385.632		DELTA=	.266		
P(3)=	4.477		DELTA=	.516		
19	407.321	.000938	16.566	.32513+00	.71051+01	-----
P(1)=	146.899		DELTA=	14.928		
P(2)=	407.321		DELTA=	.157		
P(3)=	1.898		DELTA=	.221		
20	429.751	.051496	17.422	.16323+02	.29195+00	MO
P(1)=	3371.871		DELTA=	23.320		
P(2)=	429.751		DELTA=	.025		
P(3)=	4.540		DELTA=	.049		

Figure 7. - Sample of augmented page 1 teletype output.

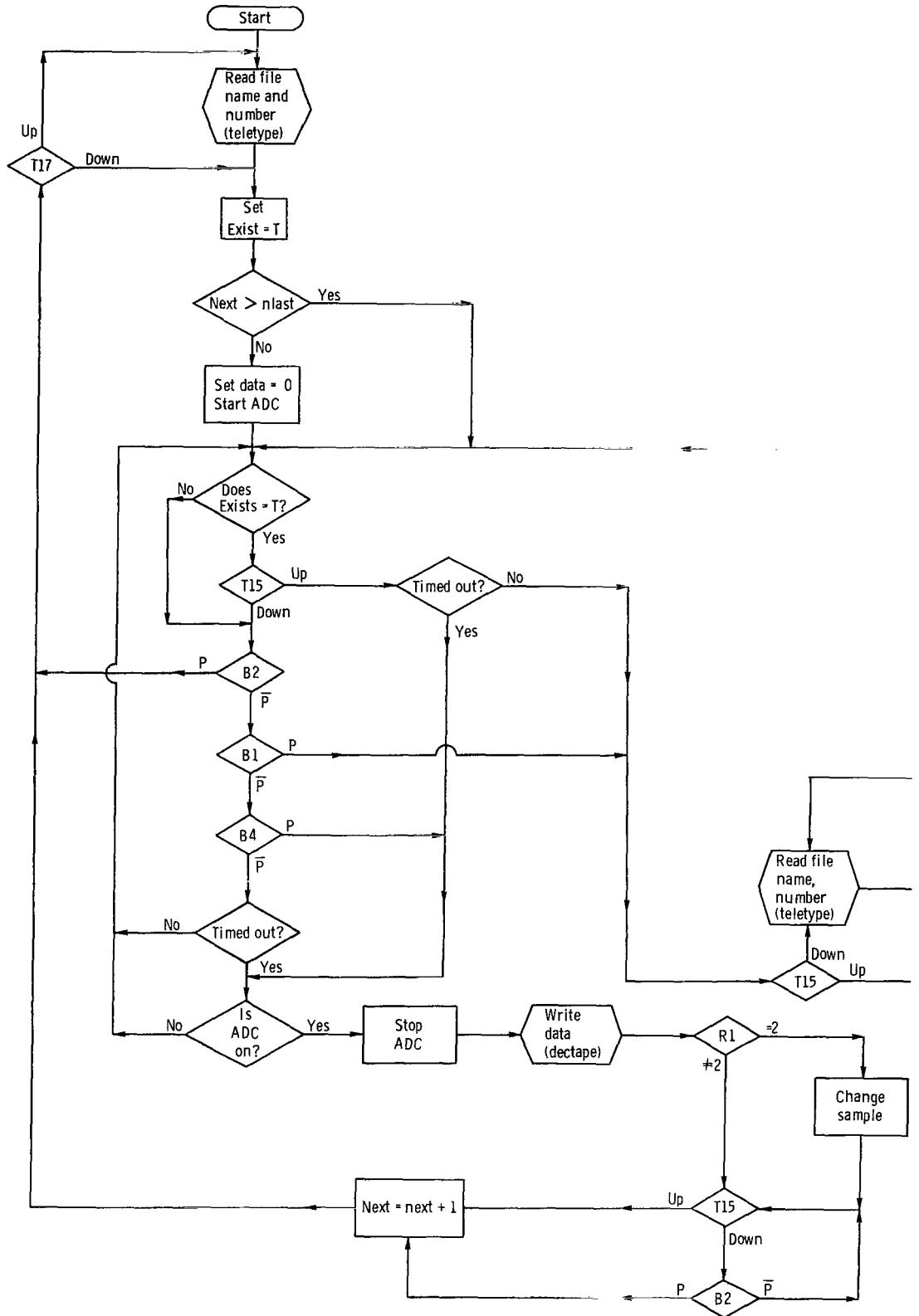


Figure 8. - Flow

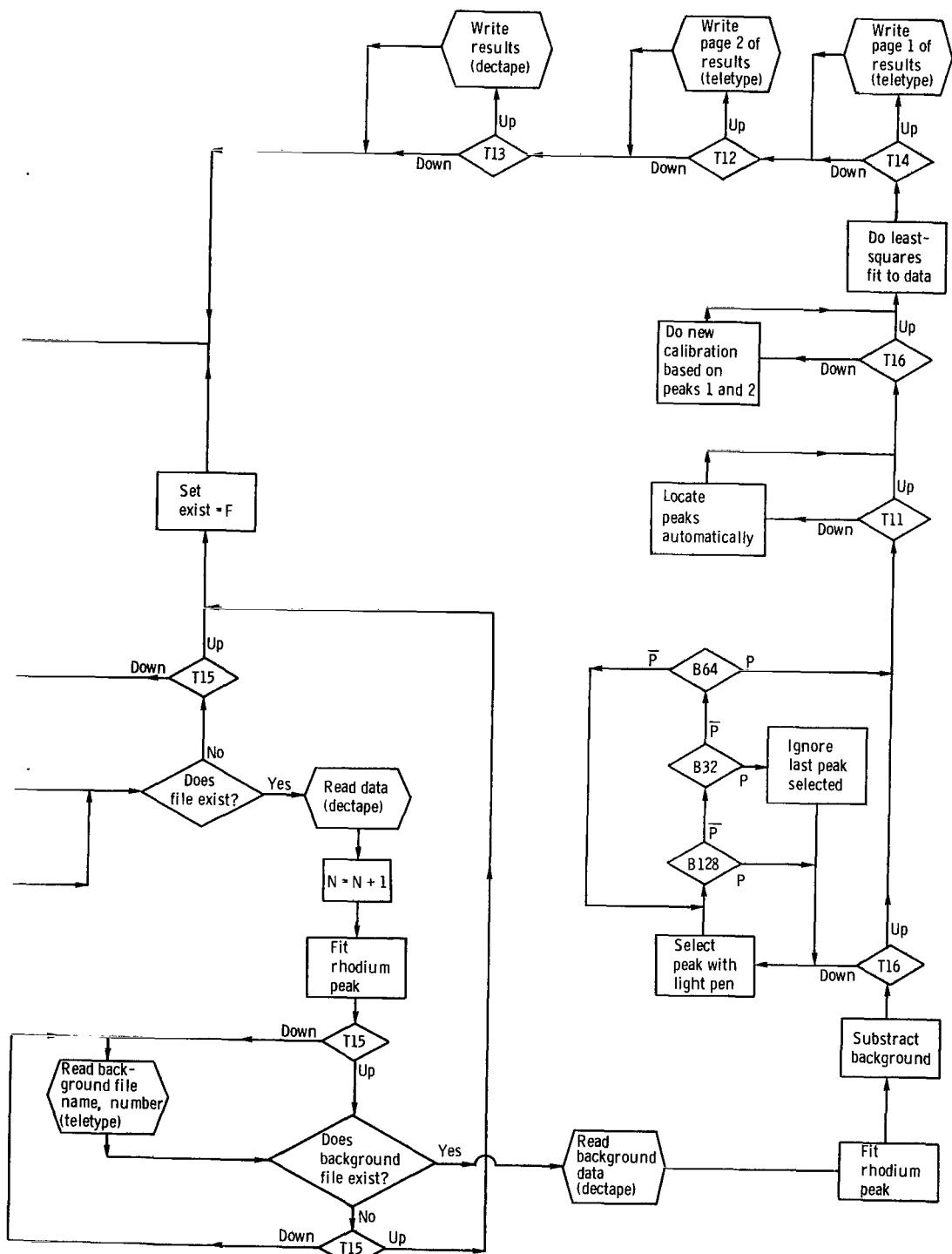


chart for program.

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